Notebook_Module1

February 20, 2023

1 Section #1.1: Linear Regression

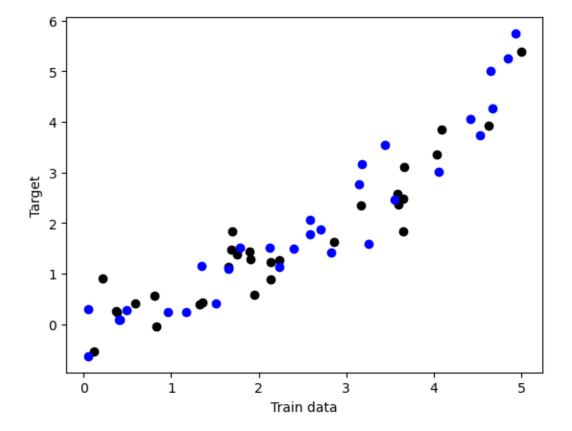
Please follow our instructions in the same order to solve the linear regression problem.

Please print out the entire results and codes when completed.

```
Train data shape: (30, 1)
Train target shape: (30,)
Test data shape: (30, 1)
Test target shape: (30,)
```

```
[]: ## PART (a):
    ## Plot the training and test data ##

plt.plot(X_train, y_train,'o', color='black')
plt.plot(X_test, y_test,'o', color='blue')
plt.xlabel('Train data')
plt.ylabel('Target')
plt.show()
```



1.1 Answer of part(a)

According to this plot, we think it is a dataset generated by a linear function plus some noises. Therefore we think a linear function could perform well.

1.2 Training Linear Regression

In the following cells, you will build a linear regression. You will implement its loss function, then subsequently train it with gradient descent. You will choose the learning rate of gradient descent to optimize its classification performance. Finally, you will get the opimal solution using closed form expression.

```
[]: from codes.Regression import Regression
```

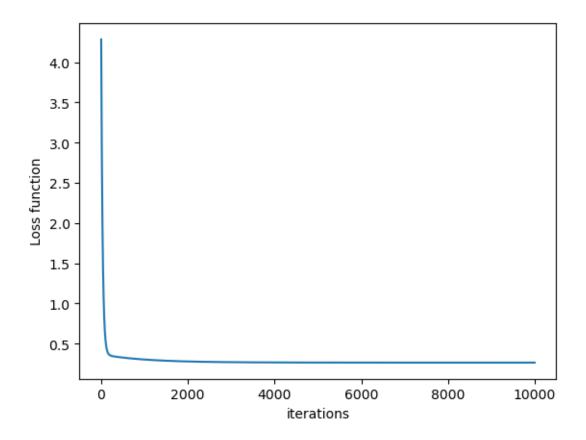
```
[]: ## PART (d):
    ## Complete loss_and_grad function in Regression.py file and test your results.
    regression = Regression(m=1, reg_param=0)
    loss, grad = regression.loss_and_grad(X_train,y_train)
    print('Loss value',loss)
    print('Gradient value',grad)
    ##
```

```
Loss value 4.282380633012861
Gradient value [[ -3.19217252]
[-10.44504584]]
```

1.3 Answer of part(d):

The loss value is about 4.28.

The gradient is vector $[-3.19 \sim, -10.44 \sim]$.



[[-0.37469204] [0.88377017]]

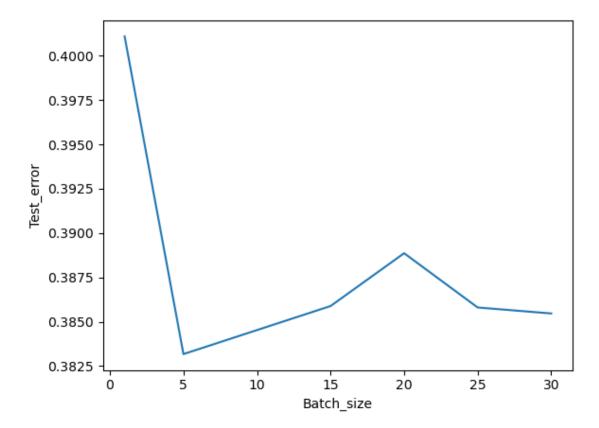
Final loss: 0.2641846537670642

```
[]: ## PART (e) (Different Batch Sizes):
    Batch = [1, 5, 15, 20, 25, 30]
    Test_err = []
    best_batch = 0
    best_test_err = 100
    # ----- #
    # YOUR CODE HERE:
    # Train the Logistic regression for different batch size Avergae the test error
     ⇔over 10 times
    # -----
    regression = Regression(m=1, reg_param=0)
    for batch in Batch:
       avg_err = 0
       for i in range(10):
           loss_history, w = regression.train_LR(X_train,y_train,_

→eta=1e-3,batch_size=batch, num_iters=10000)
           y_pred = regression.predict(X_test)
```

```
N = X_{test.shape}[0]
       y_test = np.reshape(y_test,(N,1))
       err = np.sum( np.square(y_pred - y_test) ) / N
       avg_err+=err
   test = avg_err / 10
   Test_err.append(test)
   if test < best_test_err:</pre>
       best_test_err = test
       best batch = batch
print(Test_err)
print('The best batch size is', best_batch,'the best test error is', u
 ⇒best_test_err)
learning_rate = [1e-7,1e-6,1e-5,1e-4,1e-3,1e-2,1e-1]
best_final_loss = 100
best_lr = 100
for lr in learning_rate:
   loss_history, w = regression.train_LR(X_train,y_train,_
 ⇔eta=lr,batch_size=best_batch, num_iters=10000)
   final_loss = loss_history[-1]
   if final_loss < best_final_loss:</pre>
              best_final_loss = final_loss
              best_lr = lr
print('The best learning rate for best batch size is', best_lr, 'the best final⊔
⇔loss is', best_final_loss)
# END YOUR CODE HERE
# ----- #
fig = plt.figure()
plt.plot(Batch, Test_err)
plt.xlabel('Batch_size')
plt.ylabel('Test_error')
plt.show()
fig.savefig('./plots/LR_Batch_test.pdf')
```

[0.4010980949465977, 0.38317124787831125, 0.38588178888000224, 0.38886212376210727, 0.3857999217014244, 0.38546472178456076]
The best batch size is 5 the best test error is 0.38317124787831125
The best learning rate for best batch size is 0.001 the best final loss is 0.11181510274119938



1.4 Answer of part(e)

We tested batch size ranged from [1, 5, 15, 20, 25, 30], we found the best batch size is 5 with the testing loss is 0.383.

Then we use batch size = 5, and test learning rate from [1e-7,1e-6,1e-5,1e-4,1e-3,1e-2,1e-1], then we found out the best learning rate is 0.001, with the final loss is 0.111.

```
[]: ## PART (f):
    ## Complete closed_form function in Regression.py file
    loss_2, w_2 = regression.closed_form(X_train, y_train)
    print('Optimal solution loss',loss_2)
    print('Optimal solution gradient',w_2)
```

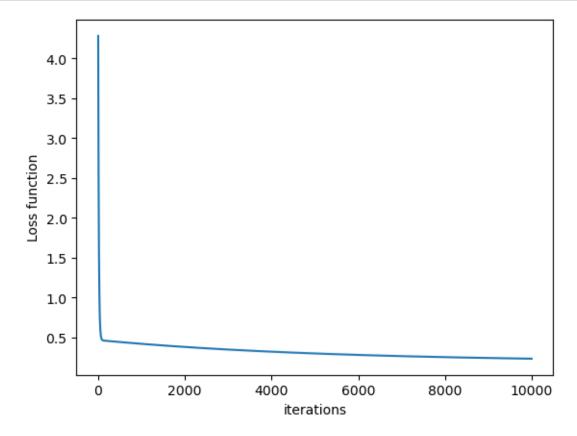
Optimal solution loss 0.26417938203964436 Optimal solution gradient [[-0.37906992] [0.8852483]]

1.5 Answer of part(f):

We get the loss and gradient of closed form: loss = 0.264~, gradient = [-0.379~, 0.885~].

The results of closed form is very close to the results we get by using numerical solution (Gradient

Descent) which is loss = 0.264~, gradient = [-0.374~, 0.884~].



```
[[0.05494868]
[0.08469512]
[0.11655802]
[0.01527028]]
Final loss when m = 3: 0.23341198948483843
```

1.6 Answer of part(g):

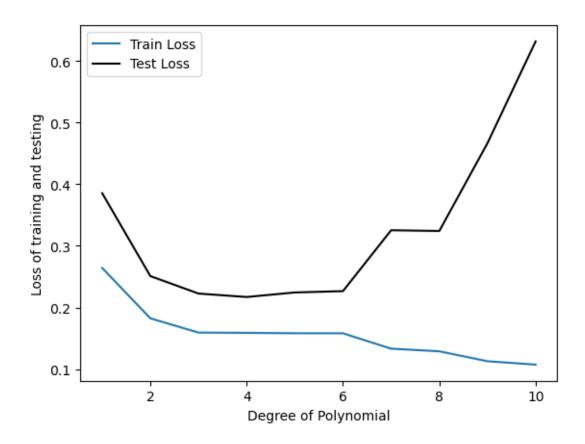
We finish the polynomial regression and test on m = 3.

The loss when m=3 is about 0.233 \sim , which is slightly better than m=1. And the gradient is $[0.0549\sim, 0.0846\sim, 0.116\sim, 0.0152\sim]$.

```
[]: ## PART (h) Overfitting
    train loss=np.zeros((10,1))
    test_loss=np.zeros((10,1))
    best_poly = -1
    best_avg_poly = -1
    best_loss = 100
    best_avg_loss = 100
    N2,d2 = X_{test.shape}
    for m in range(10):
        # print('m is', ms)
        regression = Regression(m=m+1, reg_param=0)
        loss_m, w_m = regression.closed_form(X_train, y_train)
        train_loss[m] = loss_m
        if loss_m < best_loss:</pre>
            best_loss = loss_m
            best_poly = m+1
        y_test_reshape = np.reshape(y_test, (N2,d2))
        y_test_pred = regression.predict(X_test)
        test_loss[m] = (1/N2)*np.sum(np.square(y_test_reshape-y_test_pred))
        avg = np.mean(train_loss[m] + test_loss[m])
        if avg < best_avg_loss:</pre>
            best_avg_loss = avg
            best_avg_poly = m+1
    print(train_loss)
    print( )
    print(test_loss)
    print('When degree is 3, the training loss is', train_loss[2], ', the testing_
      →loss is', test_loss[2])
    print('The best polynomial on training data is', best_poly)
    print('The best training loss of best_poly is', train_loss[best_poly-1])
    print('The best polynomial on training and testing data is', best_avg_poly)
    print('The best training loss of best_avg_poly is', train_loss[best_avg_poly-1])
    print('The best testing loss of best_avg_poly is', test_loss[best_avg_poly-1])
    y = range(1,11,1)
     # END YOUR CODE HERE
     # ----- #
    plt.plot(y, train_loss, label = 'Train Loss')
```

```
plt.plot(y, test_loss, color='black', label = 'Test Loss')
plt.xlabel('Degree of Polynomial')
plt.ylabel('Loss of training and testing')
plt.legend()
plt.show()
[[0.26417938]
 [0.18253787]
 [0.15938695]
 [0.15893654]
 [0.15829698]
 [0.15824521]
 [0.13330658]
 [0.12905253]
 [0.11292506]
 [0.1074876]]
[[0.38546511]
 [0.25101632]
 [0.22269398]
 [0.21710013]
 [0.22448356]
 [0.2266377]
 [0.32538753]
 [0.3241488]
 [0.46651551]
 [0.63156399]]
When degree is 3, the training loss is [0.15938695], the testing loss is
[0.22269398]
The best polynomial on training data is 10
The best training loss of best_poly is [0.1074876]
The best polynomial on training and testing data is 4
The best training loss of best_avg_poly is [0.15893654]
```

The best testing loss of best_avg_poly is [0.21710013]



1.7 Answer of part(h)

According to the best training loss, our best model is with degree of 10. But it is clearly that the model with degree of 10 is overfitting seriously. The plot above showed that the training loss of 10 is very low but the testing loss is very high which implies overfitting.

This plot clearly showed that with too simple model, it is underfitting with training and testing loss both too high. And with too complex model, it is overfitting with very low training loss but too high testing loss.

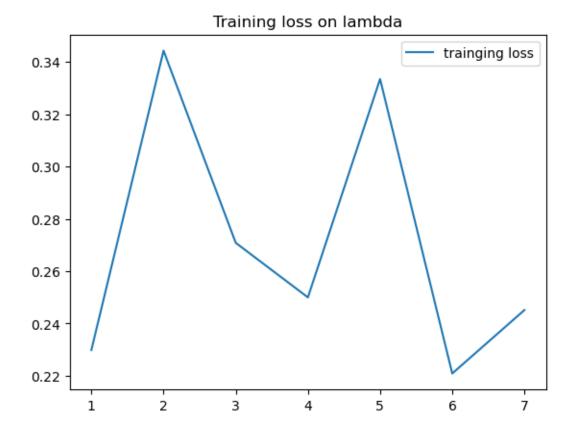
We calculated the average loss of training and testing loss and find the best degree of polynomial is 4.

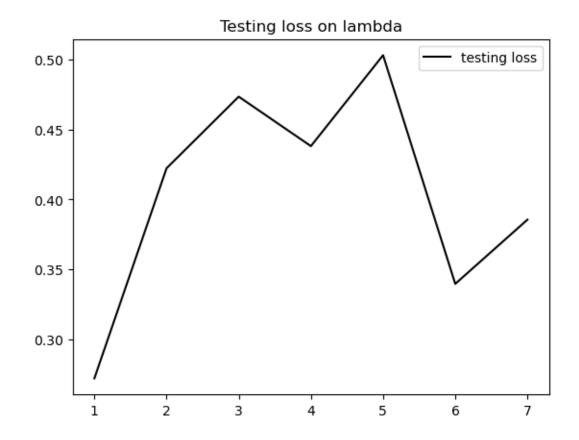
The according training loss is 0.15893654 and the according testing loss is 0.21710013.

```
# complete the following code to plot both the training and test loss in the
⇔same plot
# by taking average on 10 runs per lambda value
best_test_loss = 100
best lam = 0
for reg in range(len(lam)):
   regression = Regression(m = 3, reg param = lam[reg])
   X_aug_train = regression.gen_poly_features(X_train)
   X_aug_test = regression.gen_poly_features(X_test)
   train_err = 0
   test_err = 0
   N1,d1 = X_train.shape
   N2,d2 = X_{test.shape}
   for i in range(10):
       loss_hist, w_reg = regression.train_LR(X_train, y_train, eta=5e-4,_u
 ⇔batch_size=10, num_iters=5000)
       y_train_pred = np.dot(X_aug_train, w_reg)
       train_err+= (1/N1) * np.sum(np.square(y_train_pred - np.
 →reshape(y_train, (N1,d1))))
       y_test_reshape = np.reshape(y_test, (N2,d2))
       y_test_pred = np.dot(X_aug_test, w_reg)
       test_err+= (1/N2) * np.sum(np.square(y_test_reshape-y_test_pred))
   train err/=10
   test_err/=10
   train_loss[reg] = train_err
   test_loss[reg] = test_err
   if test err < best test loss:</pre>
       best_test_loss = test_err
       best_lam = reg
y = range(1,8,1)
print('The best lambda is', lam[best_lam], 'with best testing loss', u
⇔best_test_loss)
# ------ #
# END YOUR CODE HERE
# ------ #
plt.figure()
plt.plot(y,train_loss,label = 'trainging loss')
plt.title('Training loss on lambda')
plt.legend()
plt.show()
plt.figure()
plt.plot(y,test_loss, color='black', label = 'testing loss')
plt.legend()
```

```
plt.title('Testing loss on lambda')
plt.show()
```

The best lambda is 1e-08 with best testing loss 0.27191611411493943





1.8 Answer of part(j):

We test different regularization parameters: [1e-8, 1e-7, 1e-6, 1e-5, 1e-4, 1e-3, 1e-2]

The best param we find is 1e-8 and the testing loss is about 0.271.

2 Section #1.2: Binary Classification

Please follow our instructions in the same order to solve the binary classification problem. Please print out the entire results and codes when completed.

```
print('Train data shape: ', X_train.shape)
     print('Train target shape: ', y_train.shape)
     print('Test data shape: ',X_test.shape)
     print('Test target shape: ',y_test.shape)
    Train data shape: (5000, 784)
    Train target shape: (5000, 1)
    Test data shape: (500, 784)
    Test target shape: (500, 1)
[ ]: # PART (a):
     # To Visualize a point in the dataset
     index = 11
     X = np.array(X_train[index], dtype='uint8')
     X = X.reshape((28, 28))
     fig = plt.figure()
     plt.imshow(X, cmap='gray')
     plt.show()
     if y_train[index] == 1:
         label = 'Dress'
     else:
         label = 'Shirt'
     print('label is', label)
```

2.1 Answer of part(a)

X_train dimension: [5000, 784]X_test dimension: [500, 784]

2.2 Train Perceptron

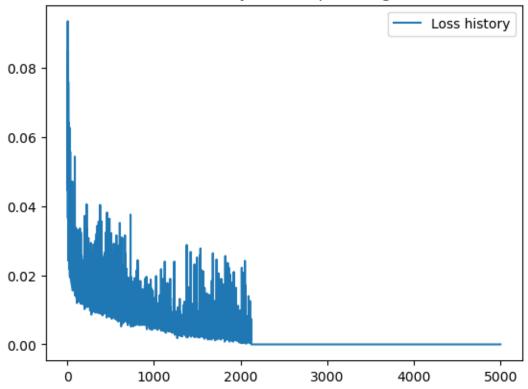
In the following cells, you will build Perceptron Algorithm.

```
# for m range from 1 to N
# ----- #
for i in range(N):
   mis_per = 0
   for j in range(N):
      train_point = X_train_h[j]
      train_label = y_train[j]
      h = np.dot(train_point, W)
      if h * train_label <= 0:</pre>
         W+= train_label * train_point.reshape(-1,1)
   y_pred = np.int_(np.sign(np.dot(X_train_h,W)))
   mis_per = np.count_nonzero(y_pred - y_train) / N
   loss_hist.append(mis_per)
print('The last value of loss history:', loss_hist[-1])
12_norm_val = (np.linalg.norm(W,ord=2)**2)
print('The squared 12 norm value of the weight:', 12_norm_val)
plt.figure()
plt.plot(loss_hist, label = 'Loss history')
plt.legend()
plt.title('The loss history of Perceptron Algorithm')
# ----- #
# END YOUR CODE HERE
# ----- #
```

The last value of loss history: 0.0
The squared 12 norm value of the weight: 659979038983.0001

[]: Text(0.5, 1.0, 'The loss history of Perceptron Algorithm')





2.3 Answer of part(b)

The last value of the loss function is 0.0

The squared norm of W is 812390.9397469916

The last misclassification is 0.0 so it could converge which implies that this dataset is linearly separable.

The percentage of misclassified points in the test data is 0.009

2.4 Answer of part(c):

The percentage of misclassification on test set is 0.9%.

2.5 Train Logistic Regression

In the following cells, you will build a logistic regression. You will implement its loss function, then subsequently train it with gradient descent.

```
[]: from codes.Logistic import Logistic
```

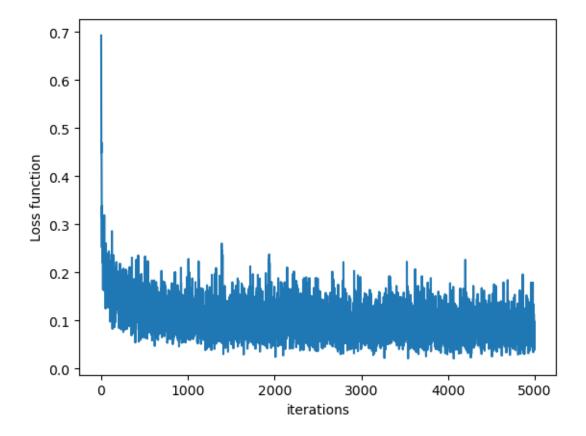
```
[]: ## PART (f):
    X_train = np.load('./data/binary_classification/X_train.npy')
    y_train = np.load('./data/binary_classification/y_train.npy')
    X_test = np.load('./data/binary_classification/X_test.npy')
    y_test = np.load('./data/binary_classification/y_test.npy')
    ## Complete loss_and_grad function in Logistic.py file and test your results.
    N,d = X_train.shape
    logistic = Logistic(d=d, reg_param=0)
    loss, grad = logistic.loss_and_grad(X_train,y_train)
    print('Loss function=',loss)
    print(np.linalg.norm(grad,ord=2)**2)
```

Loss function= 0.6931471805599454 188442.90860327004

2.6 Answer of part(f):

The loss is $0.693\sim$.

The Frobenius norm of w is about 188442.9~



Weight squared norm 0.0002852379284381107 Final loss 0.06647200567712908

2.7 Answer of part(g):

The final loss is about $0.0664\sim$.

The Frobenius norm of w is about $0.000285\sim$.

4.8 %

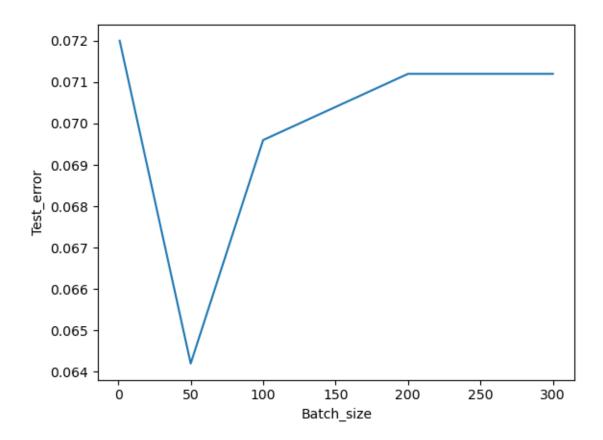
2.8 Answer of part(h):

The percentage of misclassification on test set is 4.8%.

```
[]: ## PART (i):
    Batch = [1, 50, 100, 200, 300]
    test_err = np.zeros((len(Batch),1))
    # ======= #
    # YOUR CODE HERE:
    # Train the Logistic regression for different batch size. Avergae the test_{\sqcup}
    ⇔error over 10 times
    # ----- #
    for i in range(len(Batch)):
       batch_size = Batch[i]
       test_error = 0
       num_iters = int(6000 / batch_size)
       for j in range(10):
          loss_history, w = logistic.train_LR(X_train,y_train,_
     ⇔eta=1e-5,batch_size=batch_size, num_iters=num_iters)
          y_test_pred = logistic.predict(X_test)
          test_error += np.sum((y_test!=y_test_pred))/X_test.shape[0]
          # print(test_error)
       test_error/=10
       test_err[i] = test_error
    min_idx = np.argmin(test_err)
    best_batchsize = Batch[min_idx]
    best_test_err = min(test_err)
    print('The best batch size is', best_batchsize,'the best test error is', u
     ⇒best_test_err)
    # END YOUR CODE HERE
    # ----- #
    fig = plt.figure()
    plt.plot(Batch,test_err)
    plt.xlabel('Batch_size')
    plt.ylabel('Test_error')
    plt.show()
    fig.savefig('./plots/LR_Batch_test.pdf')
```

```
/Users/yiwenzhang/Desktop/PROJECT-1/codes/Logistic.py:61: RuntimeWarning: overflow encountered in exp grad = (1/N) *((1/(1+np.exp(-h)).T - y.T)@X_aug).T

The best batch size is 50 the best test error is [0.0642]
```



2.9 Answer of part(i)

The batch size of 50 gives out the best result: testing err is 6.42%

2.10 Train SVM

In the following cells, you will build SVM. You will implement its loss function, then subsequently train it with mini-batch gradient descent. You will choose the learning rate of gradient descent to optimize its classification performance. Finally, you will get the best regularization parameter.

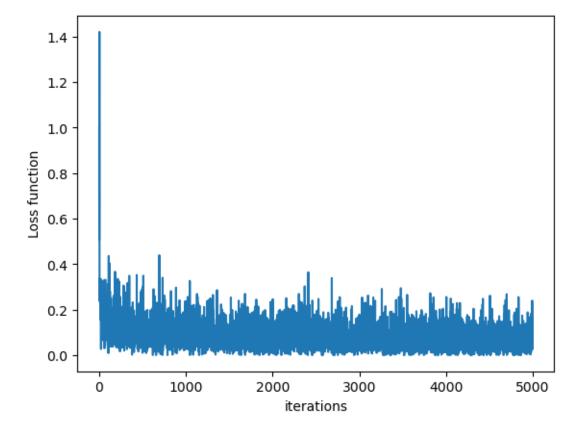
```
[]: from codes.SVM import SVM

[]: ## PART (l):
    ## Complete loss_and_grad function in Logistic.py file and test your results.
    N,d = X_train.shape
    svm = SVM(d=d, reg_param=0)
    loss, grad = svm.loss_and_grad(X_train,y_train)
    print('Loss function=',loss)
    print(np.linalg.norm(grad,ord=2)**2)
```

Loss function= 1.0 753771.6344130802

2.11 Answer of part(l):

The loss is 1.0. The Frobenius norm of w is about 753771.~



- 0.00026929379564440007
- 0.16983518000000017

2.12 Answer of part(m)

The final loss is 0.169.

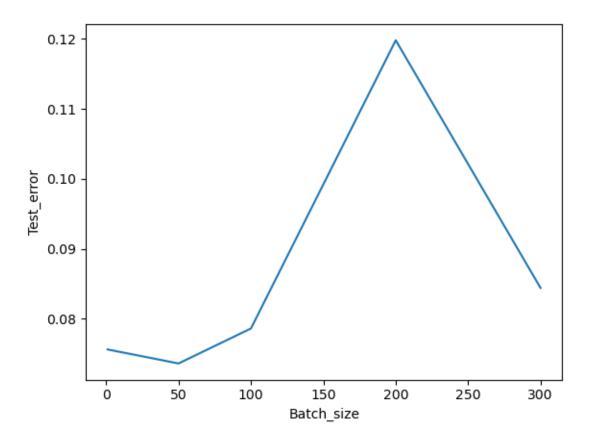
The Frobenius norm of grad is 0.000269.

6.000 %

2.13 Answer of part(n)

The percentage of misclassification on test set is 6.00%

```
[]: # PART (o)
    Batch = [1, 50, 100, 200, 300]
    test_err = np.zeros((len(Batch),1))
    # YOUR CODE HERE:
    # Train the SVM for different batch size Avergae the test error over 10 times
    for i in range(len(Batch)):
       batch_size = Batch[i]
       test_error = 0
       num_iters = int(6000 / batch_size)
       for j in range(10):
          svm = SVM(d=d, reg_param=0)
          loss_history, w = svm.train_svm(X_train,y_train,_
     ⇔eta=1e-5,batch_size=batch_size, num_iters=num_iters)
          y_test_pred = svm.predict(X_test)
          err = 1 - np.mean(svm.predict(X_test) == (y_test > 0))
          test_err[i] += 0.1*err
    fig = plt.figure()
    plt.plot(Batch,test_err)
    plt.xlabel('Batch_size')
    plt.ylabel('Test_error')
    plt.show()
    fig.savefig('./plots/svm_Batch_test.pdf')
```



The best batch size is: 50 with the best testing err, [0.0736]

2.14 Answer of part(o):

We found that the best batch size here is 50.

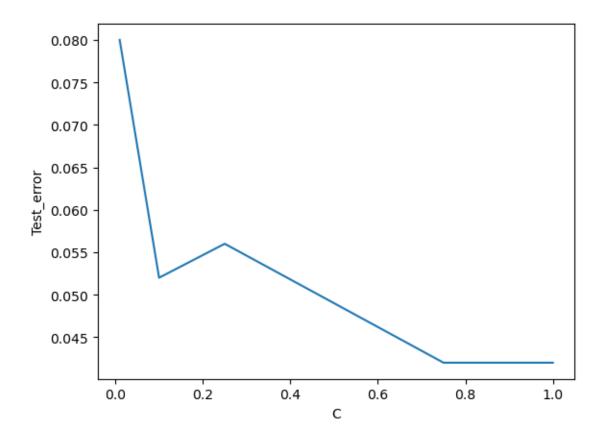
The corresponding test error is 7.36%

2.15 Train Kernelized SVM

In the following cells, you will build a RBF-SVM. You will use sklearn.svm.SVC package and its functions to train a RBF SVM. You will try to select a regularization parameter C and compare your results with non-kernelized SVM.

```
[]: # PART (p)
from sklearn.svm import SVC
```

```
[]: C_range = [0.01,0.1, 0.25, 0.75, 1] # inverse regularization parameter
   test_err = np.zeros((len(C_range),1))
   # Notes about sklearn.SVM
   # example definition of SVM classifier: clf = SVC (kernel='rbf', C = C)
   # Use clf.fit (X_train, y_train.ravel()) to train the model
   # Use 2*(clf.decision\_function(X\_test)>0)-1 to get the classification decisions\Box
    \hookrightarrow (y pred)
   # YOUR CODE HERE:
   # Train the kernelized SVM for different C
   # ----- #
   for i in range(len(C_range)):
      myc = C_range[i]
      myclf = SVC(kernel='rbf', C = myc)
      myclf.fit(X_train, y_train.ravel())
      acc = myclf.score(X_test,y_test)
      test err[i] = 1 - acc
   # END YOUR CODE HERE
   fig = plt.figure()
   plt.plot(C_range,test_err)
   plt.xlabel('C')
   plt.ylabel('Test_error')
   plt.show()
   fig.savefig('./plots/svm_C_test.pdf')
```



```
[]: best_err = min(test_err)
   idx = np.argmin(test_err)
   best_C = C_range[idx]
   print('The best C is:', best_C, 'with the best testing err,', best_err)
```

The best C is: 0.75 with the best testing err, [0.042]

2.16 Answer of part(p)

The best C for kernelized SVM is 0.75 and the according testing error is 4.2%.

The performance of kernelized SVM is better than unkernelized SVM generally.

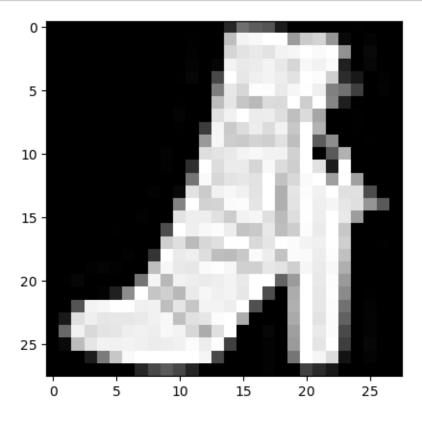
3 Section #1.3: Multi-Class Logistic Regression and Adaboost

Please follow our instructions in the same order to solve the linear regression problem. Please print out the entire results and codes when completed.

```
[]: X_train, y_train = mnist_reader.load_mnist('./data/fashion-mnist', kind='train')
X_test, y_test = mnist_reader.load_mnist('./data/fashion-mnist', kind='t10k')
print('Train data shape: ', X_train.shape)
print('Train target shape: ', y_train.shape)
```

Train data shape: (60000, 784)
Train target shape: (60000,)
Test data shape: (10000, 784)
Test target shape: (10000,)

```
# PART (a)
# To Visualize a point in the dataset
index = 11
X = np.array(X_train[index], dtype='uint8')
X = X.reshape((28, 28))
fig = plt.figure()
plt.imshow(X, cmap='gray')
plt.show()
fig.savefig('./plots/Sample.pdf')
print('label is', label[y_train[index]])
```



label is Ankle boot

3.1 Answer of part(a)

The dimension of X_train and X_test is 60000 x 784 and 10000 x 784

3.2 Train Multi-Class Logistic Regression

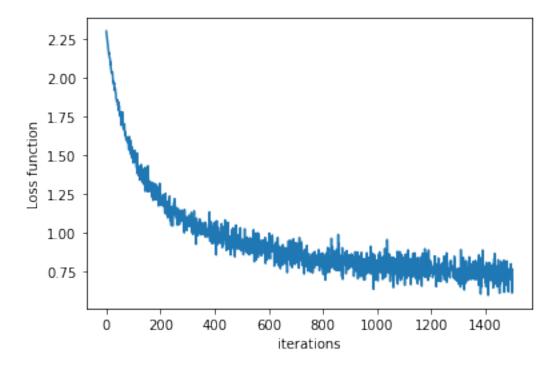
In the following cells, you will build a Multi-Class logistic regression. You will implement its loss function, then subsequently train it with gradient descent. You will implement L1 norm regularization, and choose the best regularization parameter.

```
[]: from codes.MLogistic import MLogistic
```

```
[]: ## PART (d)
    ## Complete loss_and_grad function in Logistic.py file and test your results.
    num_classes = len(np.unique(y_train))
    num_features = X_train.shape[1]

logistic = MLogistic(dim=[num_classes,num_features], reg_param=0)
loss, grad = logistic.loss_and_grad(X_train[:5000],y_train[:5000])
    print('Loss function=',loss)
    print('Frobenius norm of grad=',np.linalg.norm(grad))
##
```

Loss function= 2.3025850929940463 Frobenius norm of grad= 424.07889418407046



- 0.01122956070669866
- 0.7088758061416439

3.3 Answer of part(e)

The final value of the loss function is .709.

The final value of the square Frobenius Norm is .011.

training error: 0.23098333333333332 test error: 0.244099999999998

3.4 Answer of part(f)

Percentage of misclassfied points is 24.4%

```
[]: ## PART (h):
    reg = [0, 1e-6, 1e-3, 1e-2, 1e-1, 1]
    train_err =np.zeros((len(reg),1))
    test_err =np.zeros((len(reg),1))
    # ----- #
    # YOUR CODE HERE:
    # complete the following code to plot both the training and test loss in the
     ⇔same plot
    # for m range from 1 to 10
    \#i
    print("Part i")
    for i in range(len(reg)):
        r = reg[i]
        logistic = MLogistic(dim = [num_classes, num_features], reg_param = r)
        _, _ = logistic.train_LR(X_train, y_train, eta = 1e-7, batch_size = 200,__
     \rightarrownum_iters = 3000)
        y_train_pred = logistic.predict(X_train)
        train_err[i] = (1-np.mean(np.equal(y_train, y_train_pred)))
        y_test_pred = logistic.predict(X_test)
        test_err[i] = (1-np.mean(np.equal(y_test, y_test_pred)))
        print("For reg param "+str(reg[i])+", the test error is⊔

¬"+str(test_err[i])+".")

    best param1 = reg[np.argmin(test err)]
    print("The best parameter seems to be "+str(best_param1))
    print("Part i end")
    #1.1.
    print("Part ii")
    val_err = np.zeros((len(reg),1))
    for i in range(len(reg)):
        r = reg[i]
        perm = np.random.permutation(X_train.shape[0])
        X_train_shuffled = X_train[perm]
        y_train_shuffled = y_train[perm]
        X_folds = np.split(X_train_shuffled, 5)
        y_folds = np.split(y_train_shuffled, 5)
        flds = np.arange(5)
        for k in flds:
            bef = flds[:k]
            aft = flds[k+1:]
            fold_idx = np.concatenate((bef, aft))
```

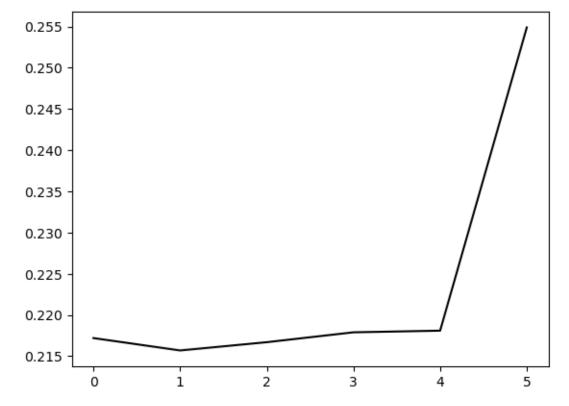
```
X_val_fold = X_folds[k]
       y_val_fold = y_folds[k]
       X_train_folds = [X_folds[j] for j in fold_idx]
       X_train_folds = np.concatenate(X_train_folds, axis = 0)
       y_train_folds = [y_folds[j] for j in fold_idx]
       y_train_folds = np.concatenate(y_train_folds, axis = 0)
       logistic = MLogistic(dim = [num_classes, num_features], reg_param = r)
        _, _ = logistic.train_LR(X_train_folds, y_train_folds, eta = 1e-7,__
 ⇔batch_size = 200, num_iters = 3000)
       y_val_pred = logistic.predict(X_val_fold)
       val_err[i] += (1-np.mean(np.equal(y_val_fold, y_val_pred)))
   val err[i] /= 5
   print("Average validation error for reg of "+str(r)+", is "+str(val_err[i]))
#best req
best_param2 = reg[np.argmin(val_err)]
print("The best parameter in this case was "+str(best_param2)+", which yields a⊔
 ⇔test error of:")
logistic = MLogistic(dim = [num_classes, num_features], reg_param = best_param2)
_, _ = logistic.train_LR(X_train, y_train, eta=1e-7, batch_size = 200,_u
 →num_iters = 3000)
y_test_pred = logistic.predict(X_test)
test_error = 1-np.mean(np.equal(y_test_pred, y_test))
print(str(test error))
print("Part ii end")
# ----- #
# END YOUR CODE HERE
# =========
fig = plt.figure()
plt.plot(test_err, color='black')
plt.show()
fig.savefig('./plots/Regularization.pdf')
y = np.arange(len(reg))
plt.figure()
plt.plot(y, val_err, label = 'Val Error with K-Folds')
plt.plot(y, test_err,color = 'black',label = 'Test Error without K-Folds')
plt.xlabel('Reg')
plt.ylabel('Error')
plt.title('Error vs Reg w/o K-Folds')
plt.legend()
plt.show()
```

Part i

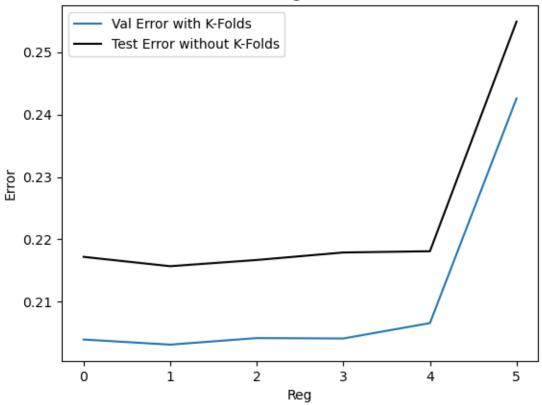
For reg param 0, the test error is [0.2172]. For reg param 1e-06, the test error is [0.2157].

```
For reg param 0.001, the test error is [0.2167].
For reg param 0.01, the test error is [0.2179].
For reg param 0.1, the test error is [0.2181].
For reg param 1, the test error is [0.2549].
The best parameter seems to be 1e-06
Part i end
Part ii
Average validation error for reg of 0, is [0.20395]
Average validation error for reg of 1e-06, is [0.20313333]
Average validation error for reg of 0.001, is [0.20418333]
Average validation error for reg of 0.01, is [0.20411667]
Average validation error for reg of 0.1, is [0.20658333]
Average validation error for reg of 1, is [0.24258333]
The best parameter in this case was 1e-06, which yields a test error of:
0.2160999999999996
Part ii end
```









3.5 Answer of part(h)

From the err plots, we could see that the test err is higher when reg is too large, for example, when reg = 1. And the test err genearly stays the same when reg is samll.

We get the best reg parameter as 1e-6 with or without using K-folds both.

We could also observe that by using K-folds, the error is generally lower: from $21\sim\%$ to $20\sim\%$. It means that our model has better generalization by using K-folds.

We retrain the model under the condition of k-folds and the best reg parameter 1e-6, and we got the test err 21.6%.

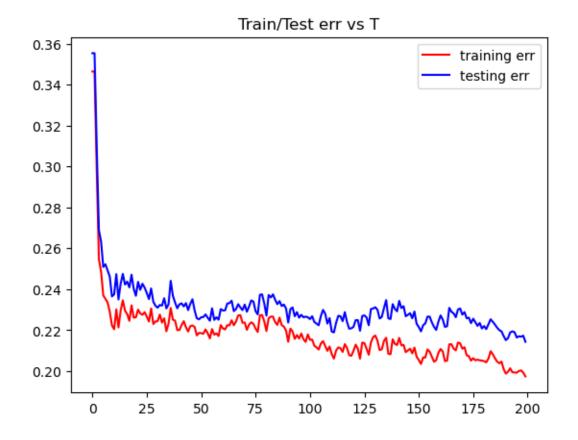
```
[]: from sklearn.tree import DecisionTreeClassifier
```

```
[]: ## PART (i):
    T = 200
    N = X_train.shape[0]
    num_classes = len(np.unique(y_train))
    num_features = X_train.shape[1]
    train_err = np.zeros((T,1))
    test_err = np.zeros((T,1))
```

```
D = np.full((N,), 1/N)
classifiers = []
coeffs = np.zeros((T,1))
for m in range(T):
   tree = DecisionTreeClassifier(max_depth=4)
   tree.fit(X_train, y_train, sample_weight=D)
   classifiers.append(tree)
   y_train_pred = tree.predict(X_train)
   alg_err = np.sum(D * (y_train != y_train_pred))
   coeffs[m] = np.log((1-alg_err)/alg_err) + np.log(num_classes-1)
   D = D*np.exp(coeffs[m]*(y_train != y_train_pred).astype(int))
   D = D/np.sum(D)
   y_train_pred = np.zeros((len(X_train), num_classes))
   for i, tree in enumerate(classifiers):
        train_pred = classifiers[i].predict(X_train)
       for k in range(num_classes):
            y_train_pred[:, k] += coeffs[i] * (train_pred == k)
   y_train_pred = np.argmax(y_train_pred, axis=1)
   y_test_pred = np.zeros((len(X_test), num_classes))
   for i, tree in enumerate(classifiers):
       test_pred = classifiers[i].predict(X_test)
       for k in range(num classes):
            y_test_pred[:, k] += coeffs[i] * (test_pred == k)
   y_test_pred = np.argmax(y_test_pred, axis=1)
   train_err[m] = (1-np.mean(np.equal(y_train_pred, y_train)))
   test_err[m] = (1-np.mean(np.equal(y_test_pred, y_test)))
    # print(train_err[m])
    # print(test_err[m])
plt.plot(train_err,color = 'red',label = 'training err')
plt.plot(test_err,color='blue',label = 'testing err')
```

```
[]: fig = plt.figure()
  plt.plot(train_err,color = 'red',label = 'training err')
  plt.plot(test_err,color='blue',label = 'testing err')
  plt.title('Train/Test err vs T')
  plt.legend()
  plt.show()
  fig.savefig('./plots/Adaboost.pdf')

  print('The test err when T = 1:', test_err[0])
  print('The test err when T = 200:', test_err[-1])
```



The test err when T = 1: [0.3554] The test err when T = 200: [0.2143]

3.6 Answer of part(i)

The test error when t = 1 is 35.54%.

The test error when t = 200 is 21.43%.

3.7 Contributions

This project is finished by 2 persons, Robert Ozturk and Yiwen Zhang, equally.

Appendix_Regression

February 20, 2023

```
[]: # %load regression.py
    import numpy as np
    class Regression(object):
       def __init__(self, m=1, reg_param=0):
           Inputs:
             - m Polynomial degree
             - regularization parameter reg_param
           Goal:
            - Initialize the weight vector self.w
            - Initialize the polynomial degree self.m
            - Initialize the regularization parameter self.reg
           11 11 11
           self.m = m
           self.reg = reg_param
           self.dim = [m+1, 1]
           self.w = np.zeros(self.dim)
       def gen_poly_features(self, X):
           Inputs:
            - X: A numpy array of shape (N,1) containing the data.
           Returns:
            - X_out an augmented training data to an mth degree feature vector e.g.
     \hookrightarrow [1, X, X<sup>2</sup>, ..., X<sup>m</sup>].
           11 11 11
           N,d = X.shape
           m = self.m
           X_out= np.zeros((N,m+1))
           if m==1:
              # ----- #
              # YOUR CODE HERE:
              # IMPLEMENT THE MATRIX X_out=[1, X]
              # ----- #
              X_out = np.hstack((np.ones((N,1)), X))
              # END YOUR CODE HERE
```

```
# ======== #
    else:
        # ----- #
        # YOUR CODE HERE:
        # IMPLEMENT THE MATRIX X_{out}=[1, X, x^2, \ldots, X^m]
        X_out = np.hstack((np.ones((N,1)), X))
        for i in range(m-1):
           X_{power} = np.zeros((N,1))
          X_power = X_power + (i + 2)
          new_X = np.power(X,X_power)
          X_out = np.hstack((X_out, new_X))
        # ----- #
        # END YOUR CODE HERE
        # ================= #
    return X_out
 def loss_and_grad(self, X, y):
     n n n
    Inputs:
     - X: N x d array of training data.
     - y: N x 1 targets
    Returns:
     - loss: a real number represents the loss
     - grad: a vector of the same dimensions as self.w containing the \Box
⇔gradient of the loss with respect to self.w
    11 11 11
    loss = 0.0
    grad = np.zeros_like(self.w)
    m = self.m
    N,d = X.shape
    if m==1:
        # ----- #
        # YOUR CODE HERE:
        # Calculate the loss function of the linear regression
        # save loss function in loss
        # Calculate the gradient and save it as grad
        # ------ #
       y_pred = self.predict(X)
       y=np.reshape(y,(N,1))
       reg = np.abs(self.w) * self.reg
        loss = (1/N) * np.sum(np.square(y - y_pred)) + np.sum(reg)
       X_out = self.gen_poly_features(X)
        dw_l1_norm = np.sign(self.w) * self.reg
        grad = (np.dot(X_out.T, y_pred - y)) * 2 / N + dw_l1_norm
```

```
\# qrad = (np.dot(X_out.T, y_pred - y)) * 2 / N
        # ----- #
        # END YOUR CODE HERE
        # ----- #
     else:
        # YOUR CODE HERE:
        # Calculate the loss function of the polynomial regression with_
⇔order m
        y_pred = self.predict(X)
       y=np.reshape(y,(N,1))
       reg = np.abs(self.w) * self.reg
       loss = (1/N) * np.sum(np.square(y - y_pred)) + np.sum(reg)
       X_out = self.gen_poly_features(X)
       dw_l1_norm = np.sign(self.w) * self.reg
       grad = (np.dot(X_out.T, y_pred - y)) * 2 / N + dw_l1_norm
        # ----- #
        # END YOUR CODE HERE
        # ------ #
    return loss, grad
 def train_LR(self, X, y, eta=1e-3, batch_size=1, num_iters=1000) :
     Finds the coefficients of a \{d-1\} \hat{} th degree polynomial
     that fits the data using least squares batch gradient descent.
     Inputs:
     - X
              -- numpy array of shape (N,1), features
              -- numpy array of shape (N,), targets
              -- float, learning rate
     - num_iters -- integer, maximum number of iterations
    Returns:
     - loss_history: vector containing the loss at each training iteration.
     - self.w: optimal weights
    loss_history = []
    N,d = X.shape
    for t in np.arange(num_iters):
       X_batch = None
       y_batch = None
        # YOUR CODE HERE:
        # Sample batch size elements from the training data for use in
⇔gradient descent.
```

```
# After sampling, X_batch should have shape: (batch_size,1),_
→y_batch should have shape: (batch_size,)
       # The indices should be randomly generated to reduce correlations_
\hookrightarrow in the dataset.
       # Use np.random.choice. It is better to user WITHOUT replacement.
       # ----- #
       num_train = N
       idx = np.random.choice(range(num_train), batch_size, replace=False)
       X_{batch} = X[idx]
       y_batch = y[idx]
       # ----- #
       # END YOUR CODE HERE
       loss = 0.0
       grad = np.zeros_like(self.w)
       # YOUR CODE HERE:
       # evaluate loss and gradient for batch data
       # save loss as loss and gradient as grad
       # update the weights self.w
       # ----- #
       loss, grad = self.loss_and_grad(X_batch,y_batch)
       self.w = self.w - eta * grad
       # END YOUR CODE HERE
       # ----- #
       loss history.append(loss)
    return loss_history, self.w
 def closed_form(self, X, y):
    11 11 11
    Inputs:
    - X: N x 1 array of training data.
    - y: N x 1 array of targets
    Returns:
    - self.w: optimal weights
    m = self.m
    loss = 0
    N,d = X.shape
    if m==1:
       # ----- #
       # YOUR CODE HERE:
       # obtain the optimal weights from the closed form solution
       # ----- #
       X_out = self.gen_poly_features(X)
       X_transpose = X_out.T
       y = np.reshape(y,(N,1))
```

```
X_T_y = np.dot(X_transpose, y)
     X_T_X_inv = np.linalg.inv(np.dot(X_transpose, X_out))
     self.w = np.dot(X_T_X_inv, X_T_y)
     y_pred = np.dot(X_out, self.w)
     loss = np.sum(np.square(y_pred - y)) / N
     # ============================ #
     # END YOUR CODE HERE
     # ------ #
  else:
     # YOUR CODE HERE:
     # IMPLEMENT THE MATRIX X_{out}=[1, X, x^2, \ldots, X^m]
     X_out = self.gen_poly_features(X)
     X_transpose = X_out.T
     y = np.reshape(y,(N,1))
     X_T_y = np.dot(X_transpose, y)
     X_T_X_inv = np.linalg.inv(np.dot(X_transpose, X_out))
     self.w = np.dot(X_T_X_inv, X_T_y)
     y_pred = np.dot(X_out, self.w)
     loss = np.sum(np.square(y_pred - y)) / N
     # ----- #
     # END YOUR CODE HERE
     # ------ #
  return loss, self.w
def predict(self, X):
   11 11 11
  Inputs:
   - X: N x 1 array of training data.
  Returns:
   - y_pred: Predicted targets for the data in X. y_pred is a 1-dimensional
    array of length N.
  y_pred = np.zeros(X.shape[0])
  m = self.m
  if m==1:
     # YOUR CODE HERE:
     # PREDICT THE TARGETS OF X
     # ----- #
     X_aug=self.gen_poly_features(X)
     # print(self.w)
     y_pred = np.dot(X_aug, self.w)
```

```
# print(y_pred)
 # END YOUR CODE HERE
 # ----- #
else:
 # YOUR CODE HERE:
 # IMPLEMENT THE MATRIX X_{out}=[1, X, x^2, \ldots, X^m]
 X_aug=self.gen_poly_features(X)
 # print(self.w)
 y_pred = np.dot(X_aug, self.w)
 # ----- #
 # END YOUR CODE HERE
 # ----- #
return y_pred
```

Appendix_Logistics

February 20, 2023

```
[]: # %load Logistic.py
   import numpy as np
   class Logistic(object):
      def __init__(self, d=784, reg_param=0):
          Inputs:
           - d: Number of features
           - regularization parameter reg_param
          Goal:
           - Initialize the weight vector self.w
          - Initialize the regularization parameter self.reg
         self.reg = reg_param
         self.dim = [d+1, 1]
         self.w = np.zeros(self.dim)
      def gen_features(self, X):
          HHH
          Inputs:
          - X: A numpy array of shape (N,d) containing the data.
          - X_out an augmented training data to a feature vector e.g. [1, X].
          n n n
         N,d = X.shape
         X_out= np.zeros((N,d+1))
          # ----- #
          # YOUR CODE HERE:
          # IMPLEMENT THE MATRIX X out=[1, X]
          # ============= #
         X_out = np.hstack((np.ones((N,1)), X))
          # ----- #
          # END YOUR CODE HERE
          return X out
      def loss_and_grad(self, X, y):
          Inputs:
```

```
- X: N x d array of training data.
     -y: N x 1 labels
     Returns:
     - loss: a real number represents the loss
     - grad: a vector of the same dimensions as self.w containing the ⊔
⇔gradient of the loss with respect to self.w
     loss = 0.0
     grad = np.zeros_like(self.w)
     # print(grad.shape)
     N,d = X.shape
     # ----- #
     # YOUR CODE HERE:
     # Calculate the loss function of the logistic regression
     # save loss function in loss
     # Calculate the gradient and save it as grad
     # ----- #
     X_aug = self.gen_features(X)
     y = (y > 0).astype('float')
     h = np.dot(X_aug, self.w)
     h_{exp} = np.exp(h) + 1
     h_{\log} = np.\log(h_{\exp})
     loss = 1/N*np.sum(h_log) - 1/N*np.sum(y*h)
     grad = (1/N) *((1/(1+np.exp(-h)).T - y.T)@X_aug).T
     # END YOUR CODE HERE
     return loss, grad
  def train_LR(self, X, y, eta=1e-3, batch_size=1, num_iters=1000) :
     Inputs:
      - X
                -- numpy array of shape (N,d), features
                -- numpy array of shape (N,), labels
      - y
                -- float, learning rate
      - num_iters -- integer, maximum number of iterations
     Returns:
      - loss_history: vector containing the loss at each training iteration.
      - self.w: optimal weights
     11 11 11
     loss_history = []
     N,d = X.shape
     for t in np.arange(num_iters):
            X_batch = None
            y_batch = None
```

```
<u>_</u>______#
          # YOUR CODE HERE:
          # Sample batch_size elements from the training data for use in_
\rightarrow gradient descent.
           # After sampling, X_batch should have shape: (batch_size,1), ___
→y_batch should have shape: (batch_size,)
          # The indices should be randomly generated to reduce__
⇔correlations in the dataset.
          # Use np.random.choice. It is better to user WITHOUT_
\hookrightarrow replacement.
          num_train = N
          idx = np.random.choice(range(num_train), batch_size,__
→replace=False)
          X_batch = X[idx]
          y_batch = y[idx]
 # END YOUR CODE HERE
_______ #
          loss = 0.0
          grad = np.zeros_like(self.w)
          #__
_____#
          # YOUR CODE HERE:
          # evaluate loss and gradient for batch data
          # save loss as loss and gradient as grad
          # update the weights self.w
          #__
 ------ #
          loss, grad = self.loss_and_grad(X_batch, y_batch)
          self.w-=eta * grad
<u>_</u>______#
          # END YOUR CODE HERE
             ------ #
          loss_history.append(loss)
    return loss_history, self.w
 def predict(self, X):
     HHHH
     Inputs:
```

Appendix_SVM

February 20, 2023

```
[]: # %load SVM.py
   import numpy as np
   class SVM(object):
      def __init__(self, d=784, reg_param=0):
          Inputs:
           - d: Number of features
           - regularization parameter reg_param
          - Initialize the weight vector self.w
          - Initialize the regularization parameter self.reg
         self.reg = reg_param
         self.dim = [d+1, 1]
         self.w = np.zeros(self.dim)
      def gen_features(self, X):
          HHH
          Inputs:
          - X: A numpy array of shape (N,d) containing the data.
          - X_out an augmented training data to a feature vector e.g. [1, X].
          n n n
         N,d = X.shape
         X_out= np.zeros((N,d+1))
          # ----- #
          # YOUR CODE HERE:
          # IMPLEMENT THE MATRIX X out=[1, X]
          # ============= #
         X_out = np.hstack((np.ones((N,1)), X))
          # ----- #
          # END YOUR CODE HERE
          return X_out
      def loss_and_grad(self, X, y):
          n n n
```

```
Inputs:
     - X: N x d array of training data.
     -y: N \times 1 labels
     Returns:
     - loss: a real number represents the loss
     - grad: a vector of the same dimensions as self.w containing the
⇔gradient of the loss with respect to self.w
     11 11 11
     loss = 0.0
     grad = np.zeros_like(self.w)
     N,d = X.shape
     # YOUR CODE HERE:
     # Calculate the loss function of the SVM regression
     # save loss function in loss
     # Calculate the gradient and save it as grad
     X_aug = self.gen_features(X)
     h_wx = np.dot(X_aug, self.w)
     res = 1 - y*h_wx
     loss = (1/N) * np.sum(np.maximum(0,res))
     grad = -(1/N)*X_aug.T@(y*(res > 0))
     # ----- #
     # END YOUR CODE HERE
     return loss, grad
 def train_svm(self, X, y, eta=1e-3, batch_size=1, num_iters=1000) :
     Inputs:
      - X
               -- numpy array of shape (N,d), features
              -- numpy array of shape (N,), labels
             -- float, learning rate
      - eta
     - num_iters -- integer, maximum number of iterations
     Returns:
     - loss_history: vector containing the loss at each training iteration.
      - self.w: optimal weights
     loss_history = []
     N, d = X.shape
     for t in np.arange(num_iters):
        X_batch = None
        v batch = None
        # YOUR CODE HERE:
```

```
# Sample batch size elements from the training data for use in
⇒gradient descent.
       # After sampling, X_batch should have shape: (batch_size,1),_
→y batch should have shape: (batch size,)
       # The indices should be randomly generated to reduce correlations \Box
\hookrightarrow in the dataset.
       # Use np.random.choice. It is better to user WITHOUT replacement.
       idx = np.random.choice(range(N), batch_size, replace=False)
       X \text{ batch} = X[idx]
       y_batch = y[idx]
       # ----- #
       # END YOUR CODE HERE
       loss = 0.0
       grad = np.zeros like(self.w)
       # YOUR CODE HERE:
       # evaluate loss and gradient for batch data
       # save loss as loss and gradient as grad
       # update the weights self.w
       # ----- #
       loss, grad = self.loss_and_grad(X_batch, y_batch)
       self.w -= eta*grad
       # ----- #
       # END YOUR CODE HERE
       # ----- #
       loss_history.append(loss)
    return loss_history, self.w
 def predict(self, X):
    Inputs:
    - X: N x d array of training data.
    Returns:
    - y_pred: Predicted labels for the data in X
    y pred = np.zeros(X.shape[0])
    # ----- #
    # YOUR CODE HERE:
    # PREDICT THE LABELS OF X
    # ----- #
    X_aug = self.gen_features(X)
    h = np.dot(X_aug, self.w)
    y_pred = (h > 0).astype('int')
    # ========== #
    # END YOUR CODE HERE
```

----- # return y_pred

Appendix_Mlogistic

February 20, 2023

```
[]: # %load Mlogistic.py
   import numpy as np
   class MLogistic(object):
      def __init__(self, dim=[10,784], reg_param=0):
          Inputs:
           - dim: dimensions of the weights [number_classes X number_features]
           - reg : Regularization type [L2,L1,L]
           - regularization parameter reg_param
          Goal:
           - Initialize the weight vector self.w
          - Initialize the regularization parameter self.reg
          11 11 11
          self.reg = reg_param
          dim[1] += 1
          self.dim = dim
          self.w = np.zeros(self.dim)
      def gen_features(self, X):
          n n n
          Inputs:
          - X: A numpy array of shape (N,d) containing the data.
          Returns:
           - X out an augmented training data to a feature vector e.g. [1, X].
          N,d = X.shape
          X \text{ out= np.zeros}((N,d+1))
          # ----- #
          # YOUR CODE HERE:
          # IMPLEMENT THE MATRIX X_out=[1, X]
          X_out = np.hstack((np.ones((N,1)), X))
          # END YOUR CODE HERE
          # ----- #
          return X_out
```

```
def loss_and_grad(self, X, y):
     Inputs:
     - X: N x d array of training data.
     -y: N x 1 labels
     Returns:
     - loss: a real number represents the loss
     - grad: a vector of the same dimensions as self.w containing the
⇔gradient of the loss with respect to self.w
     11 11 11
     loss = 0.0
     grad = np.zeros_like(self.w)
     N,d = X.shape
     # ----- #
     # YOUR CODE HERE:
     # Calculate the loss function of the logistic regression
     # save loss function in loss
     # Calculate the gradient and save it as grad
     X_aug = self.gen_features(X)
     exponential = np.exp(np.dot(X_aug, self.w.T))
     sub_y = np.dot(X_aug, self.w.T)
     #Create y one-hot matrix
     y_hot = np.zeros((y.size, y.max()+1))
     y_hot[np.arange(y.size),y] = 1
     #Total loss
     loss = 1/N * (np.sum(np.log(exponential.sum(axis = 1))) - np.trace(np.

dot(sub_y, y_hot.T)))
     #Gradient
     softmax = exponential/(exponential.sum(axis = 1))[:,None]
     grad = 1/N * (np.dot((softmax - y_hot).T, X_aug))
     #req
     grad = grad + self.reg*np.sign(self.w)
     loss = loss + self.reg * np.sum(self.w)
     # END YOUR CODE HERE
     # ----- #
     return loss, grad
  def train_LR(self, X, y, eta=1e-3, batch_size=1, num_iters=1000) :
```

```
n n n
     Inputs:
      - X
               -- numpy array of shape (N,d), features
               -- numpy array of shape (N,), labels
      - y
               -- float, learning rate
      - eta
      - num_iters -- integer, maximum number of iterations
     Returns:
      - loss_history: vector containing the loss at each training iteration.
      - self.w: optimal weights
     loss_history = []
     N,d = X.shape
     for t in np.arange(num_iters):
           X_batch = None
           y_batch = None
# YOUR CODE HERE:
           # Sample batch_size elements from the training data for use in_
\rightarrow gradient descent.
           # After sampling, X_batch should have shape: (batch_size,1),\Box
→y_batch should have shape: (batch_size,)
           # The indices should be randomly generated to reduce \square
⇔correlations in the dataset.
           # Use np.random.choice. It is better to user WITHOUT_
→replacement.
 ------ #
           num_train = N
           idx = np.random.choice(range(num_train), batch_size,__
→replace=False)
           X_{batch} = X[idx]
           y_batch = y[idx]
 # END YOUR CODE HERE
           #__
_______ #
           loss = 0.0
           grad = np.zeros_like(self.w)
           #__
_______ #
           # YOUR CODE HERE:
           # evaluate loss and gradient for batch data
           # save loss as loss and gradient as grad
           # update the weights self.w
```

```
loss, grad = self.loss_and_grad(X_batch, y_batch)
        self.w-=eta * grad
        #⊔
# END YOUR CODE HERE
        #__
loss_history.append(loss)
   return loss_history, self.w
 def predict(self, X):
   11 11 11
   Inputs:
   - X: N x d array of training data.
   Returns:
   - y pred: Predicted labelss for the data in X. y pred is a 1-dimensional
     array of length N.
   y_pred = np.zeros(X.shape[0])
   # YOUR CODE HERE:
   # PREDICT THE LABELS OF X
   # ============= #
   X_aug = self.gen_features(X)
   num = np.exp(np.dot(X_aug, self.w.T))
   softmax = num/(num.sum(axis = 1))[:,None]
   y_pred = np.argmax(softmax, axis = 1)
   # END YOUR CODE HERE
   # ------ #
   return y_pred
```